



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:22 AM UTC

PDB ID : 3MDT / pdb_00003mdt
Title : Voriconazole complex of Cytochrome P450 46A1
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Deposited on : 2010-03-30
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

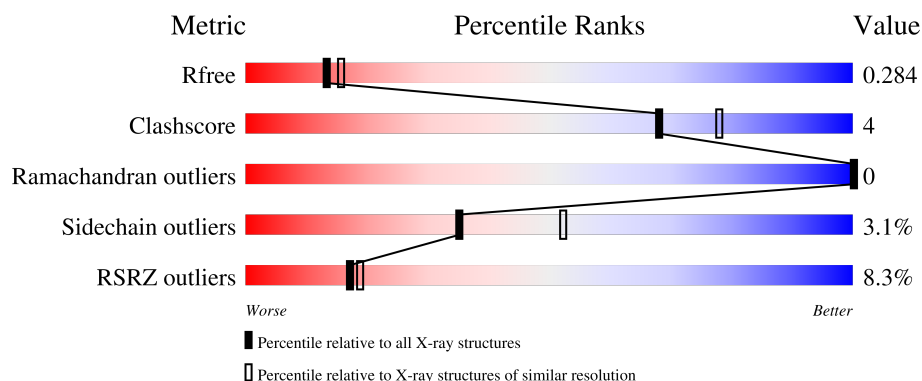
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	456	7% 83% 9% 7%
1	B	456	8% 81% 10% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	VOR	A	506	-	X	-	-
3	VOR	B	506	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7350 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Cholesterol 24-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3418	2181	598	622	17			
1	B	424	Total	C	N	O	S	0	0	0
			3418	2181	598	622	17			

There are 12 discrepancies between the modelled and reference sequences:

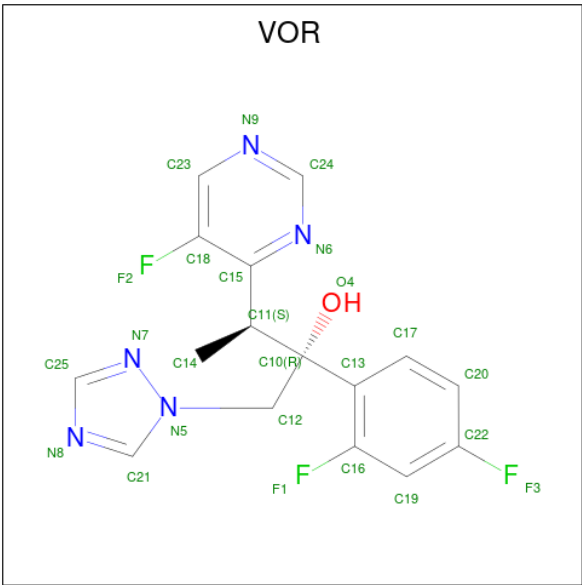
Chain	Residue	Modelled	Actual	Comment	Reference
A	49	MET	-	expression tag	UNP Q9Y6A2
A	50	ALA	-	expression tag	UNP Q9Y6A2
A	501	HIS	-	expression tag	UNP Q9Y6A2
A	502	HIS	-	expression tag	UNP Q9Y6A2
A	503	HIS	-	expression tag	UNP Q9Y6A2
A	504	HIS	-	expression tag	UNP Q9Y6A2
B	49	MET	-	expression tag	UNP Q9Y6A2
B	50	ALA	-	expression tag	UNP Q9Y6A2
B	501	HIS	-	expression tag	UNP Q9Y6A2
B	502	HIS	-	expression tag	UNP Q9Y6A2
B	503	HIS	-	expression tag	UNP Q9Y6A2
B	504	HIS	-	expression tag	UNP Q9Y6A2

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is Voriconazole (CCD ID: VOR) (formula: C₁₆H₁₄F₃N₅O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			25	16	3	5	1		
3	B	1	Total	C	F	N	O	0	0
			25	16	3	5	1		

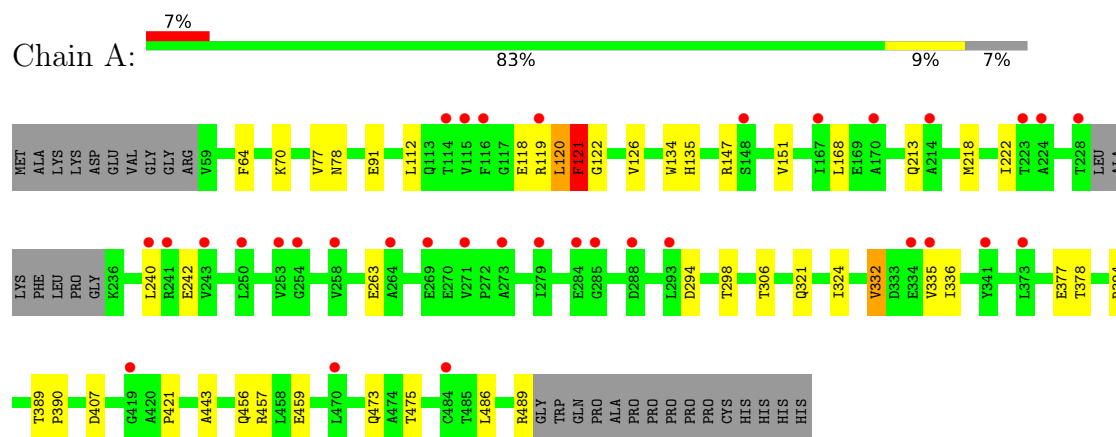
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	182	Total 182	O 182	0	0
4	B	196	Total 196	O 196	0	0

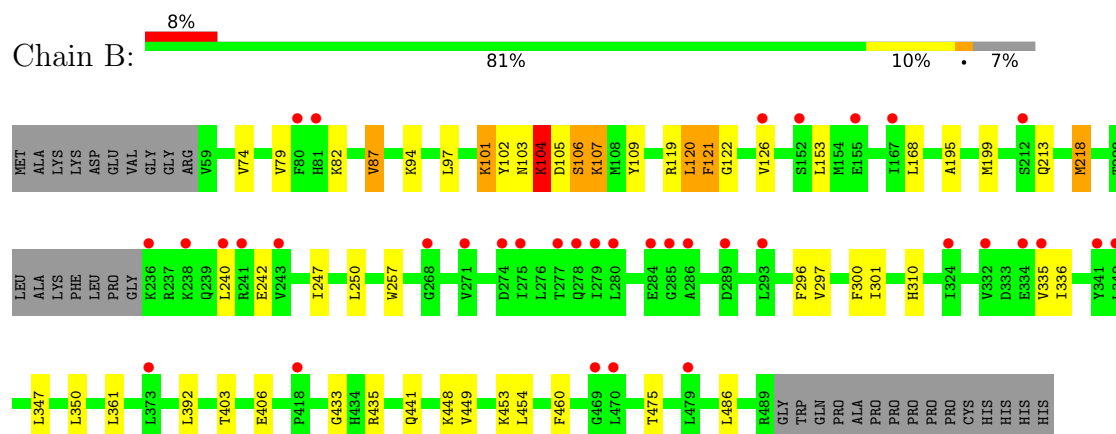
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Cholesterol 24-hydroxylase



• Molecule 1: Cholesterol 24-hydroxylase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	121.60Å 121.60Å 143.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.86 – 2.30 92.86 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.9 (92.86-2.30) 99.3 (92.86-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.229 , 0.287 0.227 , 0.284	Depositor DCC
R_{free} test set	2342 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.208 for -h,k,-l	Xtriage
Reported twinning fraction	0.500 for H, K, L 0.500 for -H, K, -L	Depositor
Outliers	1 of 46043 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7350	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, VOR

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.52	5/3485 (0.1%)	0.89	8/4700 (0.2%)
1	B	0.59	6/3485 (0.2%)	1.01	28/4700 (0.6%)
All	All	0.55	11/6970 (0.2%)	0.95	36/9400 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	120	LEU	C-N	-8.66	1.20	1.33
1	A	120	LEU	C-N	-8.52	1.20	1.33
1	B	119	ARG	C-N	-8.18	1.24	1.34
1	B	101	LYS	C-N	-7.88	1.22	1.33
1	B	106	SER	C-N	-7.21	1.24	1.33
1	A	119	ARG	NE-CZ	-6.50	1.25	1.33
1	A	119	ARG	C-N	-6.12	1.25	1.33
1	B	104	LYS	N-CA	-5.97	1.38	1.46
1	A	118	GLU	C-N	-5.85	1.25	1.33
1	A	119	ARG	CD-NE	-5.79	1.38	1.46
1	B	120	LEU	CA-C	-5.28	1.49	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	ARG	CD-NE-CZ	19.58	151.81	124.40
1	A	120	LEU	O-C-N	-11.32	107.53	122.59
1	A	119	ARG	O-C-N	-10.92	109.41	122.95
1	B	103	ASN	CA-C-N	10.69	135.89	120.82
1	B	103	ASN	C-N-CA	10.69	135.89	120.82
1	B	105	ASP	CA-C-N	9.44	133.31	120.38
1	B	105	ASP	C-N-CA	9.44	133.31	120.38
1	B	120	LEU	O-C-N	-9.06	114.12	120.83
1	B	120	LEU	CA-C-N	8.89	137.31	121.66
1	B	120	LEU	C-N-CA	8.89	137.31	121.66
1	B	119	ARG	CA-C-N	8.88	133.78	121.79
1	B	119	ARG	C-N-CA	8.88	133.78	121.79
1	A	121	PHE	O-C-N	-8.76	109.98	122.41
1	B	106	SER	CA-C-N	8.59	131.79	120.28
1	B	106	SER	C-N-CA	8.59	131.79	120.28
1	B	106	SER	CA-C-O	8.34	129.64	120.63
1	B	106	SER	O-C-N	-7.77	113.88	122.12
1	B	121	PHE	CA-C-O	7.55	128.10	119.27
1	B	103	ASN	O-C-N	7.48	132.58	123.16
1	B	105	ASP	CA-CB-CG	7.42	120.02	112.60
1	B	121	PHE	CA-CB-CG	7.40	121.20	113.80
1	A	119	ARG	CA-C-N	7.31	135.50	121.54
1	A	119	ARG	C-N-CA	7.31	135.50	121.54
1	B	107	LYS	O-C-N	7.14	129.69	122.12
1	A	121	PHE	CA-C-O	7.09	127.61	119.18
1	B	105	ASP	N-CA-C	6.74	119.11	108.67
1	B	121	PHE	O-C-N	-6.50	113.55	122.46
1	B	105	ASP	N-CA-CB	6.36	120.11	109.94
1	B	102	TYR	CA-C-N	6.24	132.92	122.87
1	B	102	TYR	C-N-CA	6.24	132.92	122.87
1	B	103	ASN	OD1-CG-ND2	5.86	128.46	122.60
1	B	120	LEU	CA-C-O	5.83	123.09	119.77
1	A	119	ARG	CA-C-O	5.67	127.39	120.92
1	B	107	LYS	CA-C-N	5.54	130.06	120.58
1	B	107	LYS	C-N-CA	5.54	130.06	120.58
1	B	120	LEU	N-CA-C	5.05	116.19	108.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	PHE	Mainchain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3418	0	3464	24	0
1	B	3418	0	3465	24	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
3	A	25	0	13	4	0
3	B	25	0	14	2	0
4	A	182	0	0	0	0
4	B	196	0	0	1	0
All	All	7350	0	7016	50	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (50) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:LEU:HD21	1:A:218:MET:HE2	1.31	1.10
1:B:120:LEU:HD11	1:B:218:MET:HE2	1.33	1.07
1:A:421:PRO:HD3	1:B:406:GLU:HB2	1.70	0.74
1:B:475:THR:HG22	3:B:506:VOR:H14	1.77	0.67
1:A:64:PHE:CE1	1:A:77:VAL:HG11	2.33	0.63
1:B:168:LEU:HB3	1:B:486:LEU:HD22	1.82	0.62
1:B:310:HIS:CE1	1:B:361:LEU:HG	2.38	0.58
1:A:77:VAL:HG12	1:A:78:ASN:N	2.21	0.55
1:B:104:LYS:HD2	1:B:435:ARG:CZ	2.41	0.50
1:A:456:GLN:HB3	1:A:457:ARG:HG2	1.93	0.50
1:B:347:LEU:HD22	1:B:448:LYS:HB3	1.93	0.50
1:A:443:ALA:HB2	2:A:505:HEM:HHC	1.94	0.50
1:A:321:GLN:HB3	1:A:324:ILE:HD12	1.93	0.50
1:B:350:LEU:HB2	1:B:448:LYS:HE2	1.94	0.50
1:A:126:VAL:HG23	1:A:298:THR:HG23	1.94	0.49
2:B:505:HEM:HMC2	2:B:505:HEM:HBC2	1.93	0.49
1:A:91:GLU:HG2	1:B:94:LYS:HD2	1.94	0.49
1:B:120:LEU:HD13	1:B:247:ILE:HG12	1.95	0.48
1:A:475:THR:HG22	3:A:506:VOR:H14	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:MET:HG3	1:B:257:TRP:CG	2.48	0.47
1:A:306:THR:OG1	3:A:506:VOR:H25	2.15	0.47
2:B:505:HEM:C1A	3:B:506:VOR:H21	2.50	0.46
1:A:147:ARG:O	1:A:151:VAL:HG23	2.16	0.46
2:A:505:HEM:HMB1	2:A:505:HEM:HBB2	1.98	0.46
1:A:122:GLY:HA3	1:A:294:ASP:CG	2.41	0.45
1:A:112:LEU:HD23	1:A:222:ILE:HD12	1.97	0.45
2:A:505:HEM:C4B	3:A:506:VOR:H25	2.51	0.45
1:B:296:PHE:O	1:B:300:PHE:HB2	2.16	0.45
1:A:218:MET:SD	1:A:218:MET:C	3.01	0.44
1:B:250:LEU:HD23	1:B:297:VAL:HG13	2.00	0.43
1:A:459:GLU:HG3	1:A:489:ARG:HG3	2.00	0.43
1:B:121:PHE:O	1:B:122:GLY:C	2.61	0.43
1:B:104:LYS:HG2	1:B:109:TYR:CE1	2.54	0.43
1:B:153:LEU:HD22	1:B:195:ALA:HB1	2.01	0.42
1:B:82:LYS:HE3	1:B:107:LYS:O	2.19	0.42
1:B:104:LYS:HE3	4:B:616:HOH:O	2.19	0.42
1:A:112:LEU:HD22	3:A:506:VOR:H23	2.01	0.42
1:B:79:VAL:HG11	1:B:392:LEU:HD12	2.01	0.42
1:A:168:LEU:HB3	1:A:486:LEU:CD2	2.50	0.42
1:A:77:VAL:CG1	1:A:78:ASN:N	2.82	0.41
1:A:389:THR:HA	1:A:390:PRO:HD3	1.90	0.41
1:B:454:LEU:HD22	1:B:460:PHE:HZ	1.84	0.41
1:A:377:GLU:OE1	1:A:384:ARG:HD3	2.21	0.41
1:A:121:PHE:O	1:A:122:GLY:C	2.61	0.41
1:B:218:MET:CE	1:B:301:ILE:HG13	2.51	0.41
1:B:74:VAL:HG22	1:B:87:VAL:HG13	2.03	0.40
1:B:97:LEU:O	1:B:433:GLY:HA3	2.21	0.40
1:B:449:VAL:O	1:B:453:LYS:HG3	2.21	0.40
1:A:134:TRP:CZ3	1:A:135:HIS:HD2	2.39	0.40
1:A:332:VAL:HG13	1:A:336:ILE:HD11	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	420/456 (92%)	407 (97%)	13 (3%)	0	100	100
1	B	420/456 (92%)	408 (97%)	12 (3%)	0	100	100
All	All	840/912 (92%)	815 (97%)	25 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	372/397 (94%)	362 (97%)	10 (3%)	39	58
1	B	372/397 (94%)	359 (96%)	13 (4%)	32	48
All	All	744/794 (94%)	721 (97%)	23 (3%)	35	52

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	LYS
1	A	213	GLN
1	A	240	LEU
1	A	242	GLU
1	A	263	GLU
1	A	332	VAL
1	A	335	VAL
1	A	378	THR
1	A	407	ASP
1	A	473	GLN
1	B	87	VAL
1	B	101	LYS
1	B	104	LYS
1	B	106	SER
1	B	126	VAL

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Mol	Chain	Res	Type
1	B	213	GLN
1	B	218	MET
1	B	240	LEU
1	B	242	GLU
1	B	335	VAL
1	B	336	ILE
1	B	403	THR
1	B	441	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	355	GLN
1	A	412	ASN
1	A	434	HIS
1	A	473	GLN
1	B	81	HIS
1	B	175	GLN
1	B	351	GLN
1	B	444	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	VOR	A	506	2	27,27,27	3.19	17 (62%)	38,39,39	3.41	24 (63%)
3	VOR	B	506	2	27,27,27	2.90	20 (74%)	38,39,39	3.12	20 (52%)
2	HEM	B	505	3,1	50,50,50	3.56	32 (64%)	67,82,82	1.82	13 (19%)
2	HEM	A	505	3,1	50,50,50	3.55	33 (66%)	67,82,82	1.69	12 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	VOR	A	506	2	-	14/21/21/21	0/3/3/3
3	VOR	B	506	2	-	14/21/21/21	0/3/3/3
2	HEM	B	505	3,1	-	6/14/54/54	-
2	HEM	A	505	3,1	-	3/14/54/54	-

All (102) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	505	HEM	FE-NB	10.17	2.26	1.94
2	B	505	HEM	FE-NB	9.98	2.25	1.94
2	B	505	HEM	C3C-C4C	-7.68	1.32	1.46
2	A	505	HEM	C3C-C4C	-7.38	1.33	1.46
2	A	505	HEM	C1C-C2C	-6.58	1.32	1.45
2	A	505	HEM	C3B-C2B	6.35	1.50	1.37
2	B	505	HEM	C1C-C2C	-6.16	1.33	1.45
2	B	505	HEM	C3B-C2B	6.11	1.49	1.37
3	A	506	VOR	C19-C16	5.79	1.47	1.37
2	B	505	HEM	C3C-C2C	5.73	1.48	1.37
2	A	505	HEM	C3C-C2C	5.73	1.48	1.37
2	B	505	HEM	CHB-C1B	5.68	1.49	1.38
2	A	505	HEM	CHB-C1B	5.68	1.49	1.38
2	A	505	HEM	CHD-C4C	5.61	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	505	HEM	CHA-C1A	5.48	1.51	1.39
2	B	505	HEM	CHC-C1C	5.46	1.49	1.38
2	B	505	HEM	CAB-C3B	-5.42	1.32	1.47
2	B	505	HEM	C1B-C2B	-5.36	1.33	1.44
2	A	505	HEM	CHC-C1C	5.30	1.48	1.38
2	A	505	HEM	C1B-C2B	-5.27	1.33	1.44
2	B	505	HEM	CHD-C4C	5.26	1.48	1.38
2	A	505	HEM	CHA-C1A	5.24	1.51	1.39
2	A	505	HEM	CAB-C3B	-5.14	1.33	1.47
3	A	506	VOR	C17-C13	5.01	1.45	1.39
3	A	506	VOR	C25-N7	4.90	1.40	1.32
3	A	506	VOR	C24-N6	4.83	1.42	1.33
2	A	505	HEM	CBB-CAB	4.46	1.51	1.30
3	B	506	VOR	C25-N7	4.45	1.40	1.32
2	B	505	HEM	CBB-CAB	4.38	1.51	1.30
3	B	506	VOR	C12-N5	-4.30	1.39	1.46
3	A	506	VOR	C23-C18	4.24	1.44	1.37
3	B	506	VOR	C13-C16	4.20	1.44	1.39
3	A	506	VOR	C18-C15	4.13	1.44	1.38
3	A	506	VOR	C15-N6	3.97	1.39	1.34
3	A	506	VOR	C14-C11	-3.97	1.47	1.53
3	B	506	VOR	C18-C15	3.79	1.44	1.38
3	A	506	VOR	C20-C17	3.75	1.44	1.38
3	B	506	VOR	C19-C22	3.65	1.43	1.37
3	B	506	VOR	C21-N5	3.62	1.40	1.33
3	B	506	VOR	C24-N6	3.54	1.40	1.33
3	B	506	VOR	C12-C10	-3.54	1.47	1.54
3	A	506	VOR	C19-C22	3.53	1.43	1.37
3	A	506	VOR	C23-N9	3.40	1.41	1.34
2	B	505	HEM	CHD-C1D	-3.38	1.31	1.39
3	A	506	VOR	C21-N5	3.35	1.39	1.33
2	B	505	HEM	C1A-NA	3.31	1.45	1.39
3	B	506	VOR	C23-C18	3.29	1.43	1.37
3	B	506	VOR	C17-C13	3.26	1.43	1.39
3	A	506	VOR	C21-N8	3.24	1.40	1.32
2	B	505	HEM	C4A-C3A	3.23	1.50	1.43
3	B	506	VOR	C19-C16	3.23	1.43	1.37
2	A	505	HEM	C4D-ND	-3.21	1.34	1.40
2	B	505	HEM	C4D-ND	-3.19	1.34	1.40
2	A	505	HEM	C4A-C3A	3.18	1.50	1.43
3	A	506	VOR	C24-N9	3.15	1.39	1.33
3	B	506	VOR	C20-C17	3.13	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	505	HEM	C1B-NB	-3.08	1.34	1.40
2	B	505	HEM	C1A-C2A	3.08	1.50	1.44
3	B	506	VOR	C21-N8	3.07	1.40	1.32
2	A	505	HEM	CHA-C4D	-3.05	1.32	1.38
2	B	505	HEM	CHA-C4D	-3.02	1.32	1.38
2	A	505	HEM	CHC-C4B	-3.01	1.32	1.39
2	B	505	HEM	FE-NC	3.01	2.05	1.95
2	A	505	HEM	CHD-C1D	-3.01	1.32	1.39
2	A	505	HEM	C1B-NB	-3.01	1.35	1.40
2	A	505	HEM	C1A-NA	2.99	1.45	1.39
2	B	505	HEM	C4A-NA	-2.97	1.34	1.39
2	B	505	HEM	C4D-C3D	2.96	1.50	1.45
2	A	505	HEM	C4A-NA	-2.95	1.34	1.39
2	A	505	HEM	C1A-C2A	2.95	1.50	1.44
3	B	506	VOR	C20-C22	2.88	1.42	1.37
2	B	505	HEM	CHC-C4B	-2.87	1.32	1.39
2	A	505	HEM	CHB-C4A	-2.84	1.33	1.39
3	A	506	VOR	C12-C10	-2.79	1.49	1.54
2	A	505	HEM	FE-NC	2.78	2.04	1.95
2	B	505	HEM	CHB-C4A	-2.76	1.33	1.39
2	B	505	HEM	C1D-ND	-2.72	1.33	1.38
3	A	506	VOR	C20-C22	2.72	1.42	1.37
3	B	506	VOR	C15-N6	2.71	1.37	1.34
2	A	505	HEM	C4D-C3D	2.63	1.49	1.45
2	B	505	HEM	C1C-NC	-2.61	1.34	1.39
3	B	506	VOR	C24-N9	2.60	1.38	1.33
2	A	505	HEM	C1C-NC	-2.60	1.34	1.39
2	A	505	HEM	C1D-ND	-2.57	1.33	1.38
3	B	506	VOR	C23-N9	2.53	1.39	1.34
2	B	505	HEM	FE-ND	2.49	2.02	1.94
2	A	505	HEM	FE-ND	2.47	2.02	1.94
2	A	505	HEM	C1D-C2D	2.43	1.49	1.44
3	B	506	VOR	C10-C11	-2.37	1.52	1.56
2	B	505	HEM	FE-NA	2.32	2.02	1.95
3	B	506	VOR	C25-N8	2.30	1.40	1.35
2	A	505	HEM	FE-NA	2.27	2.02	1.95
2	A	505	HEM	C4C-NC	-2.26	1.35	1.39
2	A	505	HEM	C3B-C4B	2.21	1.49	1.44
2	B	505	HEM	C1D-C2D	2.21	1.49	1.44
2	A	505	HEM	C4B-NB	-2.20	1.34	1.38
3	B	506	VOR	C14-C11	-2.19	1.49	1.53
2	B	505	HEM	C4B-NB	-2.14	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	506	VOR	C25-N8	2.10	1.40	1.35
2	A	505	HEM	C2A-C3A	-2.04	1.33	1.38
2	B	505	HEM	C3B-C4B	2.03	1.48	1.44
2	B	505	HEM	C2A-C3A	-2.01	1.33	1.38

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	506	VOR	C10-C12-N5	10.93	129.65	112.35
3	A	506	VOR	F1-C16-C13	9.49	129.45	119.01
3	A	506	VOR	C14-C11-C15	7.98	119.17	108.68
2	B	505	HEM	C1B-NB-C4B	5.89	112.18	105.21
3	A	506	VOR	C10-C12-N5	5.84	121.59	112.35
3	A	506	VOR	C17-C13-C10	-5.70	114.15	120.79
2	A	505	HEM	C1B-NB-C4B	5.38	111.57	105.21
2	A	505	HEM	C4B-C3B-C2B	-5.20	102.50	107.28
2	B	505	HEM	C4B-C3B-C2B	-5.15	102.55	107.28
3	B	506	VOR	C14-C11-C15	5.06	115.33	108.68
3	B	506	VOR	N9-C24-N6	-4.45	119.76	127.33
2	B	505	HEM	C3B-C4B-NB	-4.43	106.29	109.47
3	B	506	VOR	C16-C19-C22	4.39	121.35	116.67
3	B	506	VOR	F1-C16-C13	4.37	123.82	119.01
3	A	506	VOR	F1-C16-C19	-4.35	109.94	118.64
3	A	506	VOR	N9-C24-N6	-4.21	120.17	127.33
2	A	505	HEM	C3B-C4B-NB	-4.16	106.48	109.47
3	A	506	VOR	C14-C11-C10	4.16	118.65	111.55
3	A	506	VOR	C18-C15-C11	-4.01	116.78	123.14
3	B	506	VOR	C18-C15-C11	-3.97	116.84	123.14
3	A	506	VOR	C23-N9-C24	3.92	120.42	115.80
2	B	505	HEM	C4D-ND-C1D	3.87	109.79	105.21
2	A	505	HEM	C4D-ND-C1D	3.83	109.74	105.21
3	B	506	VOR	C14-C11-C10	3.82	118.07	111.55
2	B	505	HEM	CHC-C4B-NB	3.79	128.50	124.42
3	B	506	VOR	C13-C10-C11	3.78	116.19	110.88
3	A	506	VOR	C10-C13-C16	3.66	127.28	123.17
3	B	506	VOR	C23-N9-C24	3.58	120.03	115.80
2	B	505	HEM	C2A-C1A-NA	-3.54	106.23	110.15
3	B	506	VOR	C20-C22-C19	-3.54	118.56	123.23
3	B	506	VOR	C11-C15-N6	3.49	123.79	118.67
3	A	506	VOR	C10-C11-C15	-3.38	107.06	112.67
3	A	506	VOR	C11-C15-N6	3.08	123.18	118.67
3	A	506	VOR	F2-C18-C15	3.07	122.56	120.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	505	HEM	CHD-C4C-NC	3.03	127.75	124.45
3	A	506	VOR	O4-C10-C11	-3.02	104.90	110.65
3	A	506	VOR	C19-C16-C13	-3.01	120.61	123.88
2	A	505	HEM	C2A-C1A-NA	-2.99	106.83	110.15
3	A	506	VOR	C16-C19-C22	2.90	119.76	116.67
3	B	506	VOR	C25-N7-N5	2.88	109.25	102.60
2	B	505	HEM	CHD-C4C-C3C	-2.87	120.37	125.21
2	A	505	HEM	C3C-C2C-C1C	-2.81	104.39	107.05
3	A	506	VOR	C25-N7-N5	2.81	109.09	102.60
3	B	506	VOR	O4-C10-C13	-2.75	102.17	107.54
2	A	505	HEM	CHD-C4C-NC	2.75	127.45	124.45
3	B	506	VOR	C24-N6-C15	2.66	122.18	118.26
2	B	505	HEM	CHA-C1A-NA	2.65	128.66	123.86
3	A	506	VOR	N7-C25-N8	-2.65	107.35	114.90
3	B	506	VOR	N7-C25-N8	-2.62	107.42	114.90
3	B	506	VOR	C19-C16-C13	-2.59	121.06	123.88
2	B	505	HEM	O2A-CGA-CBA	2.56	122.08	114.00
2	A	505	HEM	CHC-C4B-NB	2.51	127.13	124.42
3	B	506	VOR	F2-C18-C15	2.45	122.07	120.11
3	B	506	VOR	O4-C10-C11	-2.42	106.03	110.65
2	B	505	HEM	C4C-C3C-C2C	-2.41	104.72	106.81
3	A	506	VOR	F3-C22-C19	2.40	121.70	118.28
3	A	506	VOR	O4-C10-C12	2.39	112.65	107.56
3	A	506	VOR	C12-N5-N7	2.35	124.67	121.25
3	A	506	VOR	C18-C15-N6	2.30	119.91	118.09
3	A	506	VOR	C20-C22-C19	-2.29	120.20	123.23
3	B	506	VOR	C12-N5-N7	2.29	124.59	121.25
2	A	505	HEM	CHA-C1A-NA	2.26	127.97	123.86
3	B	506	VOR	O4-C10-C12	2.23	112.30	107.56
2	B	505	HEM	C3D-C4D-ND	-2.20	107.76	110.17
3	A	506	VOR	C12-C10-C11	2.14	114.16	110.14
2	A	505	HEM	O2A-CGA-CBA	2.13	120.73	114.00
2	B	505	HEM	O2D-CGD-CBD	2.05	120.48	114.00
2	A	505	HEM	O2D-CGD-CBD	2.03	120.42	114.00
2	A	505	HEM	CHA-C4D-ND	2.02	126.86	124.37

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	505	HEM	C2B-C3B-CAB-CBB
3	A	506	VOR	O4-C10-C11-C15

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Mol	Chain	Res	Type	Atoms
3	A	506	VOR	C12-C10-C11-C14
3	A	506	VOR	C13-C10-C11-C14
3	A	506	VOR	C13-C10-C11-C15
3	A	506	VOR	O4-C10-C12-N5
3	A	506	VOR	C11-C10-C12-N5
3	A	506	VOR	C13-C10-C12-N5
3	A	506	VOR	C11-C10-C13-C16
3	A	506	VOR	C11-C10-C13-C17
3	A	506	VOR	C12-C10-C13-C16
3	A	506	VOR	C12-C10-C13-C17
3	B	506	VOR	O4-C10-C11-C15
3	B	506	VOR	C12-C10-C11-C14
3	B	506	VOR	C13-C10-C11-C14
3	B	506	VOR	C13-C10-C11-C15
3	B	506	VOR	C13-C10-C12-N5
3	B	506	VOR	C11-C10-C13-C16
3	B	506	VOR	C11-C10-C13-C17
3	B	506	VOR	C12-C10-C13-C16
3	B	506	VOR	C12-C10-C13-C17
3	A	506	VOR	O4-C10-C13-C17
3	B	506	VOR	O4-C10-C13-C17
3	A	506	VOR	O4-C10-C13-C16
3	B	506	VOR	O4-C10-C13-C16
2	A	505	HEM	C2C-C3C-CAC-CBC
3	B	506	VOR	C11-C10-C12-N5
3	A	506	VOR	C10-C12-N5-N7
3	B	506	VOR	C10-C12-N5-N7
2	B	505	HEM	C2D-C3D-CAD-CBD
2	A	505	HEM	CAD-CBD-CGD-O1D
2	B	505	HEM	CAD-CBD-CGD-O1D
2	B	505	HEM	CAD-CBD-CGD-O2D
2	A	505	HEM	CAD-CBD-CGD-O2D
2	B	505	HEM	C4B-C3B-CAB-CBB
3	B	506	VOR	O4-C10-C12-N5
2	B	505	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

4 monomers are involved in 9 short contacts:

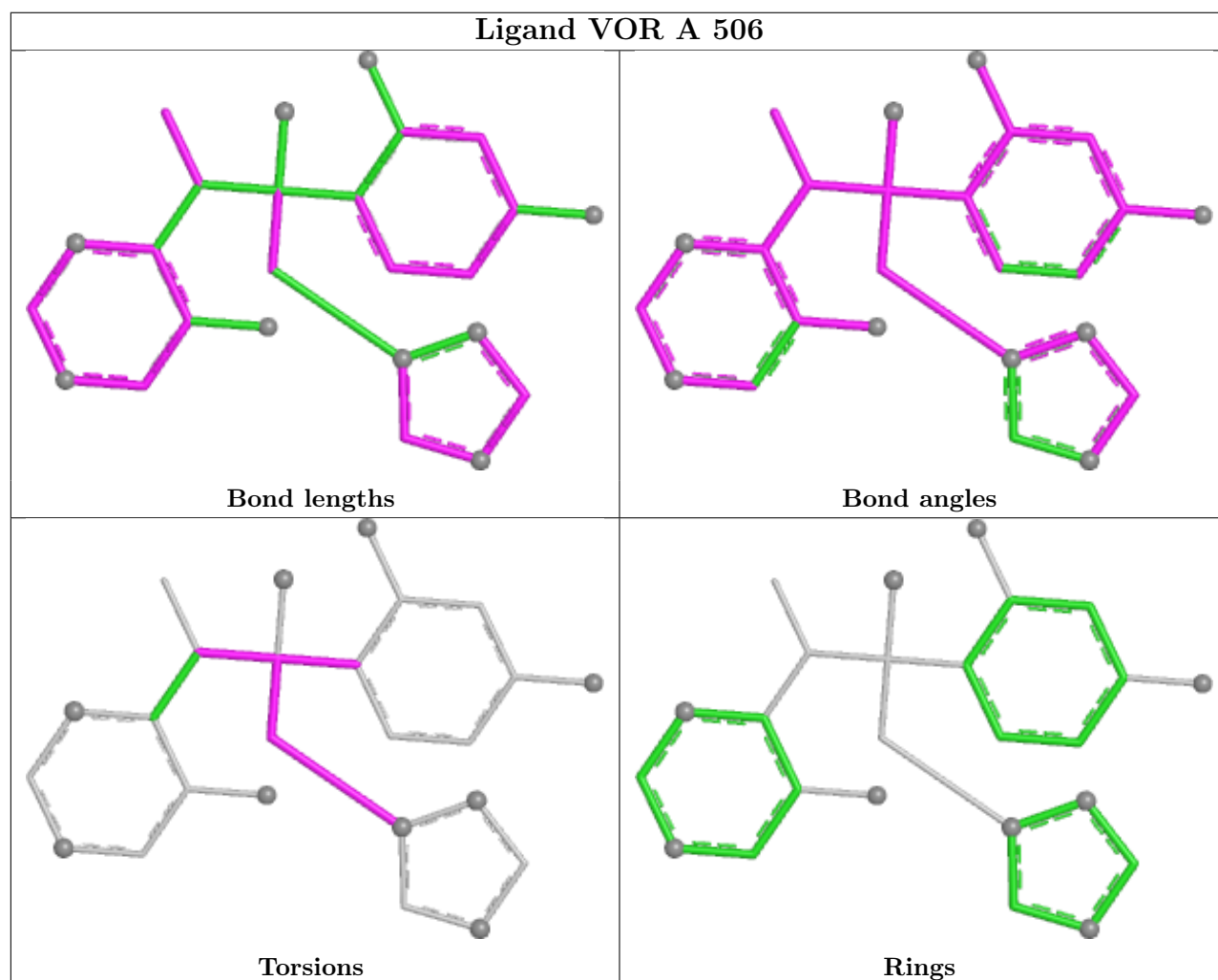
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	506	VOR	4	0
3	B	506	VOR	2	0

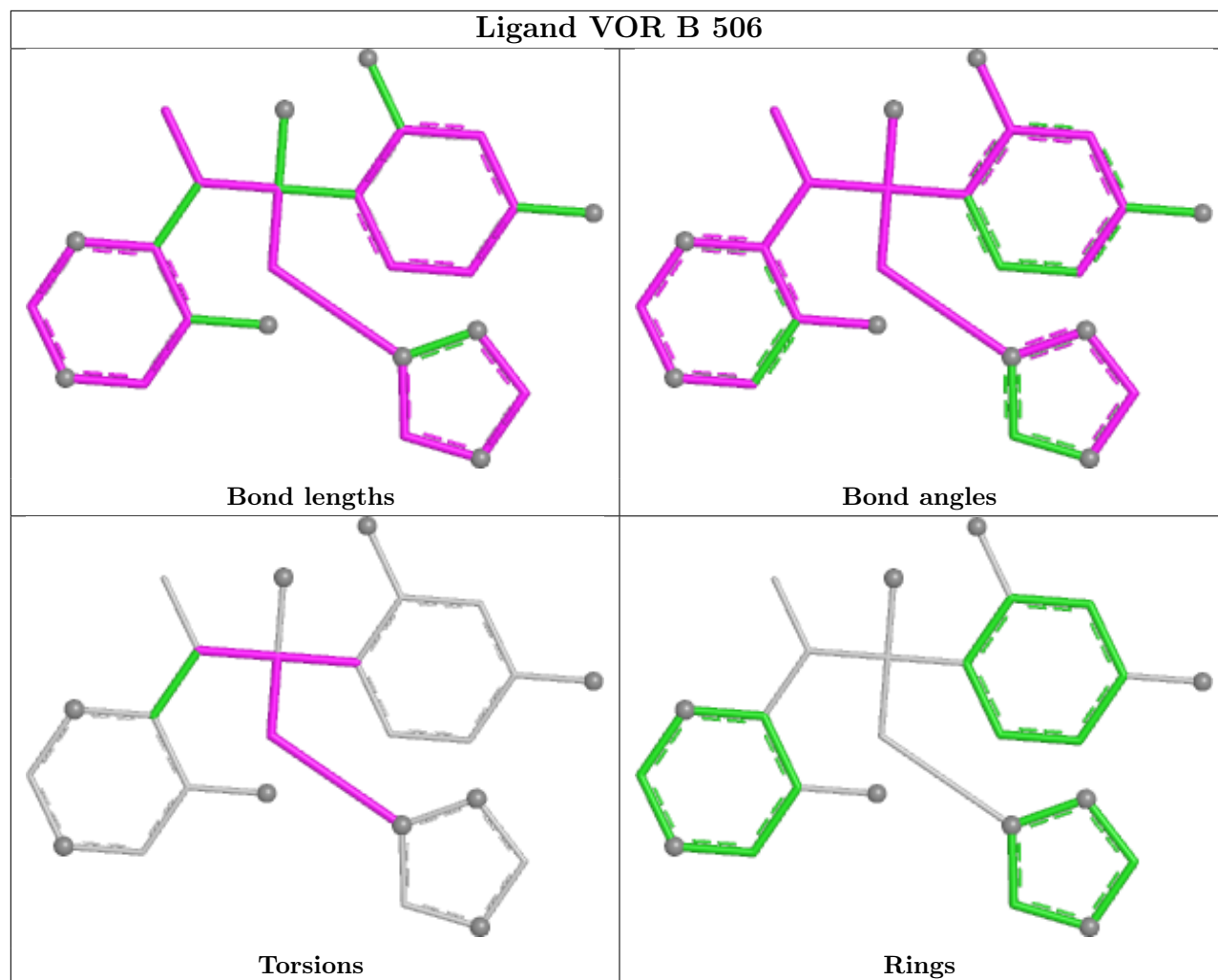
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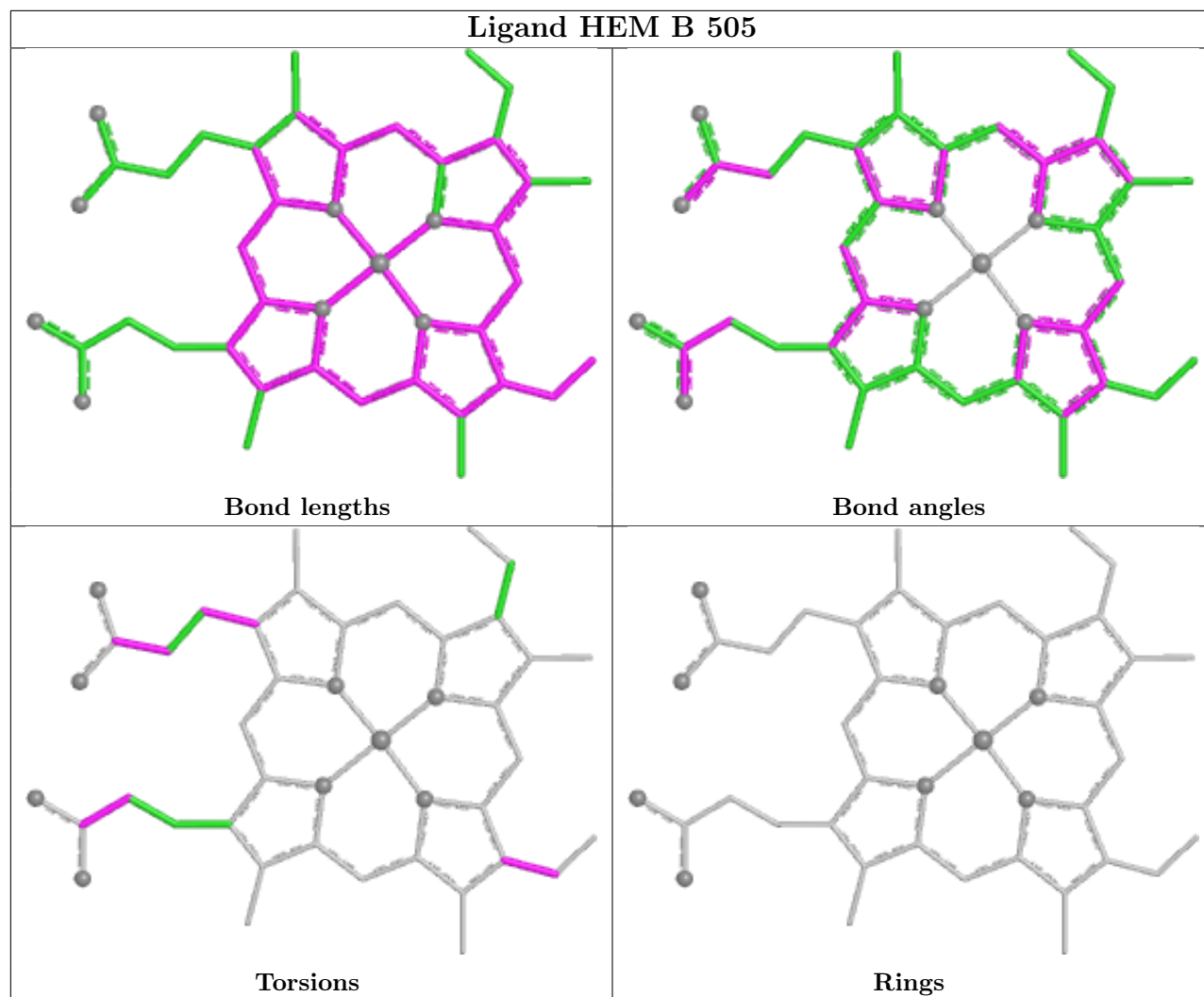
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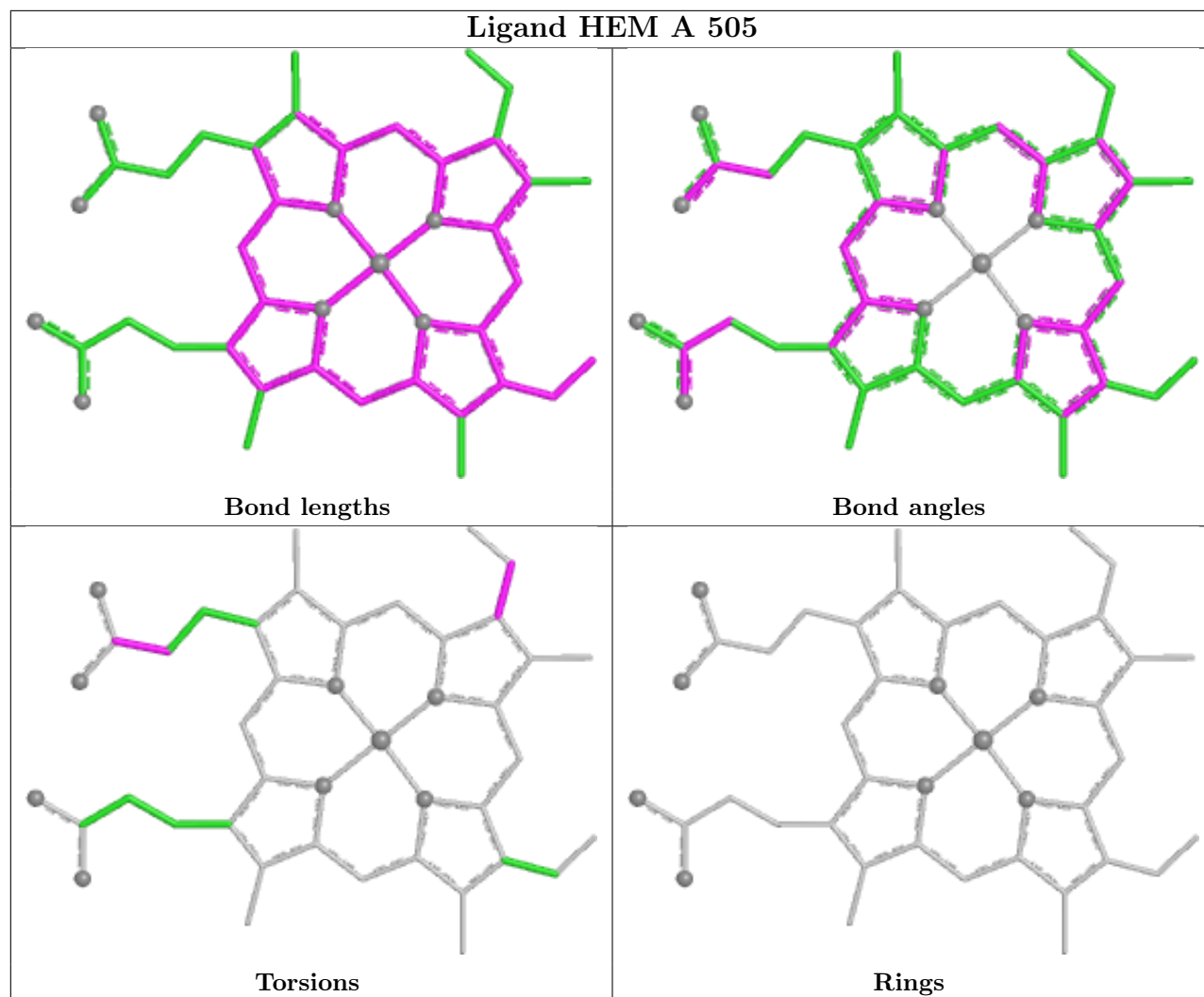
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	505	HEM	2	0
2	A	505	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	120:LEU	C	121:PHE	N	1.20

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	424/456 (92%)	0.80	34 (8%) 18 20	9, 20, 29, 35	0
1	B	424/456 (92%)	0.91	36 (8%) 16 18	8, 21, 33, 37	0
All	All	848/912 (92%)	0.85	70 (8%) 17 19	8, 20, 32, 37	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	335	VAL	4.8
1	A	224	ALA	4.5
1	B	286	ALA	3.4
1	B	289	ASP	3.4
1	B	271	VAL	3.3
1	B	277	THR	3.3
1	B	241	ARG	3.2
1	B	332	VAL	3.2
1	A	228	THR	3.2
1	B	279	ILE	3.1
1	B	418	PRO	3.0
1	A	273	ALA	3.0
1	B	275	ILE	3.0
1	A	335	VAL	2.9
1	B	126	VAL	2.9
1	A	279	ILE	2.9
1	B	469	GLY	2.9
1	A	115	VAL	2.9
1	B	167	ILE	2.9
1	A	250	LEU	2.8
1	A	148	SER	2.8
1	A	167	ILE	2.7
1	A	170	ALA	2.7
1	A	293	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	254	GLY	2.6
1	A	114	THR	2.6
1	B	341	TYR	2.6
1	B	342	LEU	2.6
1	A	341	TYR	2.6
1	B	334	GLU	2.5
1	A	419	GLY	2.5
1	B	285	GLY	2.5
1	A	116	PHE	2.5
1	B	268	GLY	2.5
1	A	269	GLU	2.5
1	B	280	LEU	2.5
1	B	373	LEU	2.5
1	A	223	THR	2.4
1	B	212	SER	2.4
1	A	214	ALA	2.4
1	B	80	PHE	2.4
1	A	288	ASP	2.4
1	A	253	VAL	2.4
1	A	240	LEU	2.4
1	B	152	SER	2.4
1	B	278	GLN	2.4
1	B	324	ILE	2.3
1	B	236	LYS	2.3
1	A	264	ALA	2.3
1	B	238	LYS	2.3
1	A	258	VAL	2.3
1	A	271	VAL	2.3
1	B	243	VAL	2.3
1	B	479	LEU	2.2
1	B	284	GLU	2.2
1	B	240	LEU	2.2
1	B	293	LEU	2.2
1	A	284	GLU	2.1
1	B	274	ASP	2.1
1	A	285	GLY	2.1
1	A	241	ARG	2.1
1	A	334	GLU	2.1
1	A	470	LEU	2.1
1	B	155	GLU	2.1
1	A	484	CYS	2.1
1	A	243	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	470	LEU	2.0
1	A	119	ARG	2.0
1	B	81	HIS	2.0
1	A	373	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

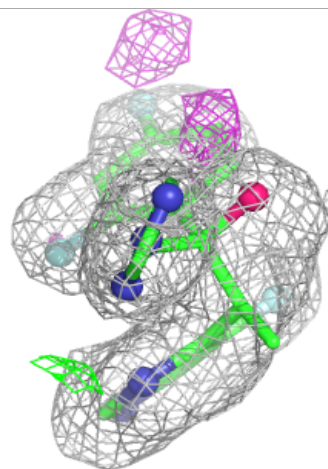
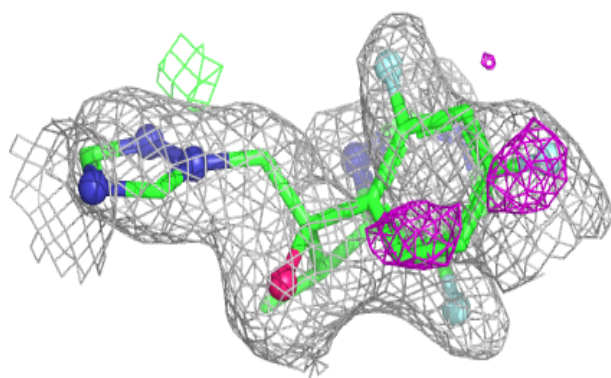
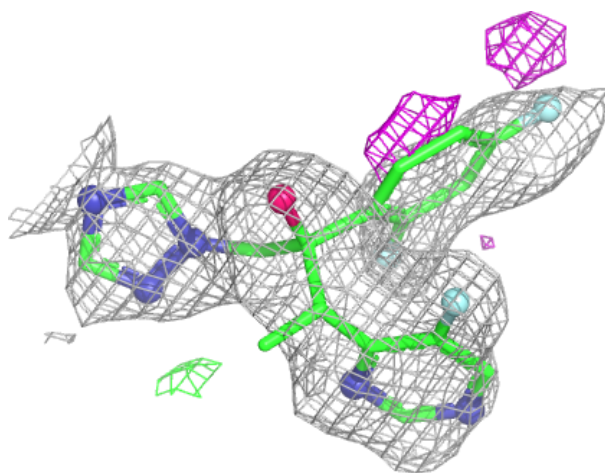
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	VOR	A	506	25/25	0.80	0.17	40,44,45,46	0
3	VOR	B	506	25/25	0.83	0.17	40,44,45,46	0
2	HEM	B	505	43/43	0.94	0.11	18,20,24,27	0
2	HEM	A	505	43/43	0.95	0.09	12,17,23,25	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

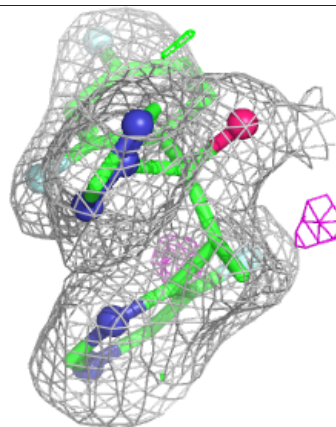
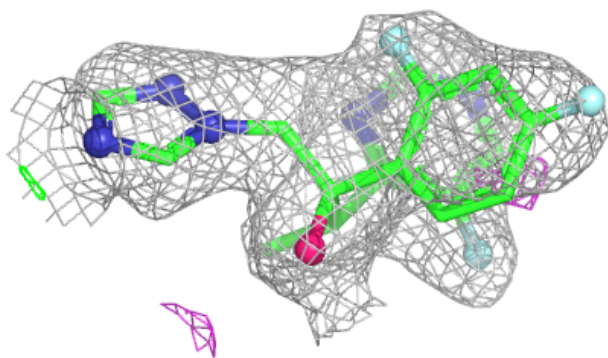
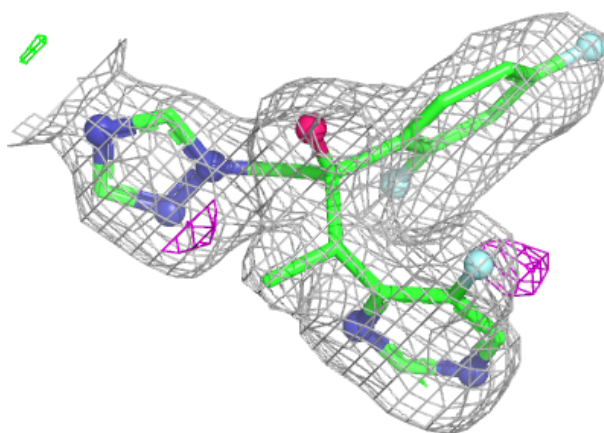
Electron density around VOR A 506:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



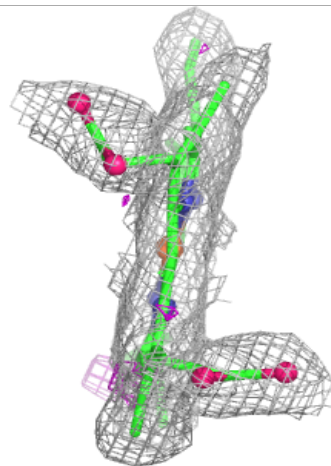
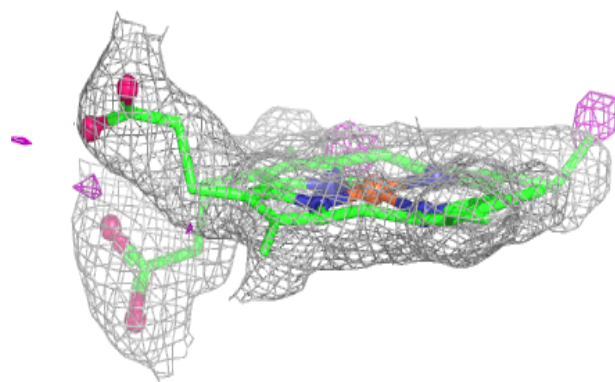
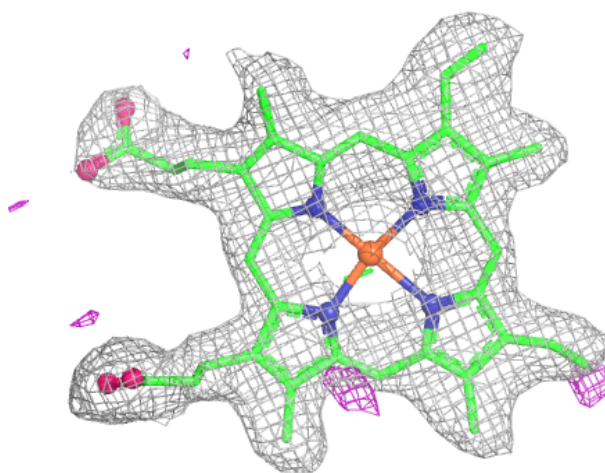
Electron density around VOR B 506:

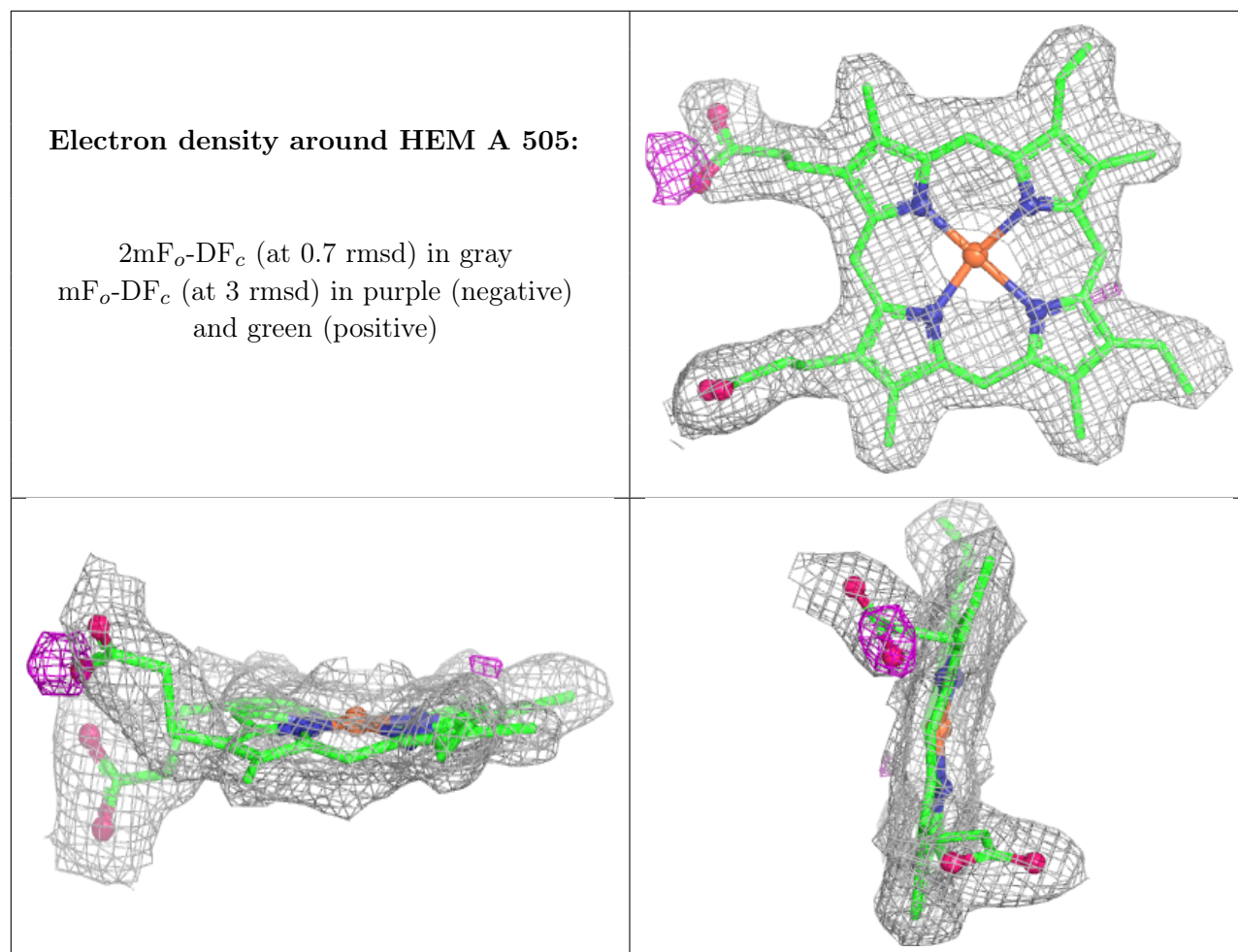
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 505:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.